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LAB-6: Parallel Cellular Algorithms and Programs

CODE:

```
#pcap

import numpy as np

# Define the problem: A simple optimization function (e.g., Sphere Function)
def optimization_function(position):
    """Example: Sphere Function for minimization."""
    return sum(x**2 for x in position)

# Initialize Parameters

GRID_SIZE = (10, 10) # Grid size (rows, columns)
NEIGHBORHOOD_RADIUS = 1 # Moore neighborhood radius
DIMENSIONS = 2 # Number of dimensions in the solution space
ITERATIONS = 30 # Number of iterations

# Initialize Population
def initialize_population(grid_size, dimensions):
    """Initialize a grid with random positions."""
    population = np.random.uniform(-10, 10, size=(grid_size[0], grid_size[1], dimensions))
    return population

# Evaluate Fitness
def evaluate_fitness(population):
    """Calculate the fitness of all cells."""
    fitness = np.zeros((population.shape[0], population.shape[1]))
    for i in range(population.shape[0]):
        for j in range(population.shape[1]):
            fitness[i, j] = optimization_function(population[i, j])
```

```
return fitness
```

```
# Get Neighborhood
```

```
def get_neighborhood(grid, x, y, radius):  
    """Get the neighbors of a cell within the specified radius."""  
    neighbors = []  
    for i in range(-radius, radius + 1):  
        for j in range(-radius, radius + 1):  
            if i == 0 and j == 0:  
                continue # Skip the current cell  
            ni, nj = x + i, y + j  
            if 0 <= ni < grid.shape[0] and 0 <= nj < grid.shape[1]:  
                neighbors.append((ni, nj))  
    return neighbors
```

```
# Update States
```

```
def update_states(population, fitness):  
    """Update the state of each cell based on its neighbors."""  
    new_population = np.copy(population)  
    for i in range(population.shape[0]):  
        for j in range(population.shape[1]):  
            neighbors = get_neighborhood(population, i, j, NEIGHBORHOOD_RADIUS)  
            best_neighbor = population[i, j]  
            best_fitness = fitness[i, j]  
  
            # Find the best position among neighbors  
            for ni, nj in neighbors:  
                if fitness[ni, nj] < best_fitness:  
                    best_fitness = fitness[ni, nj]  
                    best_neighbor = population[ni, nj]
```

```
    # Update the cell state (move towards the best neighbor)
    new_population[i, j] = (population[i, j] + best_neighbor) / 2 # Average position
return new_population
```

```
# Main Algorithm
```

```
def parallel_cellular_algorithm():
```

```
    """Implementation of the Parallel Cellular Algorithm."""
```

```
    population = initialize_population(GRID_SIZE, DIMENSIONS)
```

```
    best_solution = None
```

```
    best_fitness = float('inf')
```

```
    for iteration in range(ITERATIONS):
```

```
        # Evaluate fitness
```

```
        fitness = evaluate_fitness(population)
```

```
        # Track the best solution
```

```
        min_fitness = np.min(fitness)
```

```
        if min_fitness < best_fitness:
```

```
            best_fitness = min_fitness
```

```
            best_solution = population[np.unravel_index(np.argmin(fitness), fitness.shape)]
```

```
        # Update states based on neighbors
```

```
        population = update_states(population, fitness)
```

```
        # Print progress
```

```
        print(f"Iteration {iteration + 1}: Best Fitness = {best_fitness}")
```

```
print("\nBest Solution Found:")
```

```
print(f"Position: {best_solution}, Fitness: {best_fitness}")
```

```
# Run the algorithm
```

```
if __name__ == "__main__":  
    parallel_cellular_algorithm()
```

OUTPUT:

```
Iteration 1: Best Fitness = 0.43918427791098213  
Iteration 2: Best Fitness = 0.43918427791098213  
Iteration 3: Best Fitness = 0.062221279350329436  
Iteration 4: Best Fitness = 0.030149522005462108  
Iteration 5: Best Fitness = 0.015791278460696168  
Iteration 6: Best Fitness = 0.0025499667118763104  
Iteration 7: Best Fitness = 0.0025499667118763104  
Iteration 8: Best Fitness = 0.00019007166980743008  
Iteration 9: Best Fitness = 0.00019007166980743008  
Iteration 10: Best Fitness = 1.0432171933623911e-05  
Iteration 11: Best Fitness = 8.406928148912647e-06  
Iteration 12: Best Fitness = 5.511032710180021e-07  
Iteration 13: Best Fitness = 4.3084388056725156e-07  
Iteration 14: Best Fitness = 2.315054420755622e-07  
Iteration 15: Best Fitness = 5.245753459404661e-08  
Iteration 16: Best Fitness = 5.245753459404661e-08  
Iteration 17: Best Fitness = 4.341357920017173e-08  
Iteration 18: Best Fitness = 1.145644119860328e-08  
Iteration 19: Best Fitness = 3.147791691706415e-09  
Iteration 20: Best Fitness = 2.8192306881167533e-09  
Iteration 21: Best Fitness = 9.788374665398935e-11  
Iteration 22: Best Fitness = 9.788374665398935e-11  
Iteration 23: Best Fitness = 9.788374665398935e-11  
Iteration 24: Best Fitness = 9.788374665398935e-11  
Iteration 25: Best Fitness = 7.537171686605552e-11  
Iteration 26: Best Fitness = 7.234639306921671e-11  
Iteration 27: Best Fitness = 7.028872029493468e-11  
Iteration 28: Best Fitness = 3.340290444524624e-11  
Iteration 29: Best Fitness = 1.4953679944431498e-11  
Iteration 30: Best Fitness = 1.0817118995466254e-11  
  
Best Solution Found:  
Position: [-2.92599538e-06 -1.50188883e-06], Fitness: 1.0817118995466254e-11
```